Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE051916\
Data File : BE091839.D
Acq On : 20 May 2016 16:06
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: May 20 17:23:27 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE051916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu May 19 15:36:49 2016
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	5.000	5.000	0.0	140	0.00
2		1,4-Dioxane	0.400	0.370	7.5	111	0.00
3		n-Nitrosodimethylamine	0.400	0.364	9.0	139	0.00
4	S	2-Fluorophenol	0.400	0.419	-4.7	156	0.00
5	S	Phenol-d6	0.400	0.425	-6.2	166	0.02
6		bis(2-Chloroethyl)ether	0.400	0.389	2.8	149	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	141	0.00
8	S	Nitrobenzene-d5	0.400	0.409	-2.2	162	0.00
9		Nitrobenzene	0.400	0.395	1.3	161	0.00
10		Naphthalene	0.400	0.384	4.0	141	0.00
11		Hexachlorobutadiene	0.400	0.396	1.0	146	-0.02
12		2-Methylnaphthalene	0.400	0.369	7.8	138	-0.01
13	I	Acenaphthene-d10	5.000	5.000	0.0	122	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.378	5.5	133	0.00
15	S	2-Fluorobiphenyl	0.400	0.399	0.3	129	-0.01
16		Acenaphthylene	0.400	0.405	-1.3	130	0.00
17		Acenaphthene	0.400	0.391	2.3	127	-0.01
18		Fluorene	0.400	0.368	8.0	120	-0.01
19	I	Phenanthrene-d10	5.000	5.000	0.0	119	-0.01
20		4-Bromophenyl-phenylether	0.400	0.392	2.0	119	0.00
21		Hexachlorobenzene	0.400	0.373	6.8	112	0.00
22		Pentachlorophenol	0.400	0.253	36.8#	83	0.00
23		Phenanthrene	0.400	0.349	12.8	108	0.00
24		Anthracene	0.400	0.374	6.5	115	-0.01
25		Fluoranthene	0.400	0.377	5.8	112	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	94	-0.02
27		Benzidine	0.400	0.655	-63.7#	175	0.00
28		Pyrene	0.400	0.443	-10.7	110	0.00
29	S	Terphenyl-d14	0.400	0.436	-9.0	110	-0.02
30		Benzo(a)anthracene	0.400	0.403	-0.8	102	-0.02
31		3,3'-Dichlorobenzidine	0.400	0.557	-39.3#	150	-0.02
32		Chrysene	0.400	0.478	-19.5	105	0.00
33		Bis(2-ethylhexyl)phthalate	0.400	0.671	-67.8#	168	-0.02
34		Indeno(1,2,3-cd)pyrene	0.400	0.440	-10.0	110	-0.02
35	I	Perylene-d12	5.000	5.000	0.0	101	-0.02
36		Benzo(b)fluoranthene	0.400	0.397	0.8	106	-0.02
37		Benzo(k)fluoranthene	0.400	0.408	-2.0	108	-0.01
38	C	Benzo(a)pyrene	0.400	0.412	-3.0	110	-0.02
39		Dibenzo(a,h)anthracene	0.400	0.410	-2.5	110	-0.05
40		Benzo(g,h,i)perylene	0.400	0.403	-0.8	107	-0.05

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Compound	Amount	Calc.	%Dev	Area%	Dev(min)

(#) = Out of Range SPCC's out = 0 CCC's out = 0